Application No. 10/599,121 Docket No.: 20517/0205421-US0

Amendment dated December 29, 2009 Reply to Office Action of July 10, 2009

### AMENDMENTS TO THE CLAIMS

The following listing of the claims replaces all prior versions of the claims presented in the application.

Claim 1 (Currently amended): A method of treating a mammal suffering from or-being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a compound of the general formula (I)

wherein

 $R^1, R^2, R^3$ , and  $R^4$  independently are selected from hydrogen, optionally substituted  $C_{1-6}$ -alkyl, optionally substituted  $C_{2-6}$ -alkenyl, hydroxy, optionally substituted  $C_{1-6}$ -alkoxy, optionally substituted  $C_{1-6}$ -alkoxy, optionally substituted  $C_{1-6}$ -alkoxycarbonyl, optionally substituted  $C_{1-6}$ -alkoxycarbonyl, optionally substituted  $C_{1-6}$ -alkylcarbonyloxy, formyl, amino, monoand  $di(C_{1-6}$ -alkyl)amino, carbamoyl, mono- and  $di(C_{1-6}$ -alkyl)aminocarbonyl,  $C_{1-6}$ -alkylsulphonylamino,  $C_{1-6}$ -alkylsulphonylamino, cyano, carbamido, mono- and  $di(C_{1-6}$ -alkyl)aminocarbonylamino,  $C_{1-6}$ -alkylsulphonyl,  $C_{1-6}$ -alkylsulphonyl,  $C_{1-6}$ -alkylsulphinyl, aminosulfonyl, mono- and  $di(C_{1-6}$ -alkyl)aminosulfonyl, nitro, optionally substituted  $C_{1-6}$ -alkylthio, aryl, aryloxy, arylcarbonyl, arylamino, heterocyclyl, heterocyclyloxy, heterocyclylamino, heterocyclylcarbonyl, heteroaryl, heteroaryloxy, heteroarylamino, heteroarylcarbonyl, and halogen, where any  $C_{1-6}$ -alkyl as an amino substituent is optionally substituted with hydroxy,  $C_{1-6}$ -alkoxy, amino, mono- and  $di(C_{1-6}$ -alkyl)amino, carboxy,  $C_{1-6}$ -alkylcarbonylamino,  $C_{1-6}$ -alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted;

or  $\mathbb{R}^1$  and  $\mathbb{R}^2$  together with the carbon atoms to which they are attached form a ring;

with the proviso that R<sup>1</sup>, R<sup>2</sup>, R<sup>3</sup> and R<sup>4</sup> are not all hydrogen;

 $X^1 \text{ and } X^2 \text{ are independently selected from halogen, hydroxy} \underbrace{(-OH) \text{ and acctoxy}}_{\text{-OAc}} \underbrace{(-OAc)}_{\text{-optionally substituted}} C_{1-6}\text{-alkylearbonyloxy, amino, mono—and } \text{di}(C_{1-6}\text{-alkylearbonyloxy, amino, mono—and } \text{di}(C_{1-6}\text{-alkylearbonyloxy, meno—and } \text{di}(C_{1-6}\text{-alkylearbonylamino,} C_{1-6}\text{-alkylearbonyloxy, mercapto, optionally substituted} C_{1-6}\text{-alkylthio,} C_{1-6}\text{-alkylearbonyloxy, mercapto,} \text{optionally substituted} C_{1-6}\text{-alkylthio,} C_{1-6}\text{-alkylearbonyloxy, arylamino,} \text{heteroeyelyloxy,} \text{heteroeyelylamino, heteroaryloxy-and heteroarylamino, where any } C_{1-6}\text{-alkyl-as an amino-or-sulphur substitutent is optionally substituted with hydroxy,} C_{1-6}\text{-alkoxy, amino, mono—and } \text{di}(C_{1-6}\text{-alkylearbonylamino,} C_{1-6}\text{-alkylearbonyl, or-halogen(s),} \text{ and } \text{wherein any aryl, heteroeyelyl-and-heteroaryl-may-be-optionally-substituted;} \text{ and}$ 

pharmaceutically acceptable salts and prodrugs thereof.

Claims 2-3 (canceled).

Claim 4 (Previously presented): The method according to claim 1, wherein  $\mathbb{R}^1$  is selected from hydrogen, halogen,  $\mathbb{C}_{1\text{-}6}$ -alkyl, trifluoromethyl and  $\mathbb{C}_{1\text{-}6}$ -alkoxy.

Claim 5 (Previously presented): The method according to claim 1, wherein R<sup>2</sup> is selected from hydrogen, halogen, optionally substituted aryl, optionally substituted aryloxy, and optionally substituted heteroaryl.

Claim 6 (Previously presented): The method according to claim 1, wherein  $\mathbb{R}^3$  is selected from hydrogen, optionally substituted  $\mathbb{C}_{1.6}$ -alkoxy, halogen, cyano, optionally substituted aryl, optionally substituted aryloxy, optionally substituted heteroaryl, amino,  $\mathbb{C}_{1.6}$ -alkylcarbonylamino,  $\mathbb{C}_{1.6}$ -alkylsulphonylamino, and mono- and di( $\mathbb{C}_{1.6}$ -alkyl)aminosulfonyl.

Claim 7 (Previously presented): The method according to claim 1, wherein R<sup>4</sup> is hydrogen.

Claims 8-20 (canceled).

Claim 21 (Previously presented): The method according to claim 1, wherein R<sup>1</sup> is selected from fluoro, chloro, bromo, C<sub>1-4</sub>-alkyl, trifluoromethyl, C<sub>1-4</sub>-alkoxy, and dimethylaminocarbonyl.

Claim 22 (canceled).

Claim 23 (Previously presented): The method according to claim 1, wherein R<sup>1</sup> is selected from halogen, C<sub>1-4</sub>-alkyl, trifluoromethyl, C<sub>1-4</sub>-alkoxy, and dimethylaminocarbonyl, R<sup>2</sup> is selected from hydrogen and halogen, and R<sup>3</sup> is selected from hydrogen, halogen, C<sub>1-4</sub>-alkyl, and amino; where R<sup>2</sup> and R<sup>3</sup> are not both hydrogen.

Claim 24 (Currently amended): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IIa)

wherein

R1 is selected from hydrogen, halogen, C1-6-alkyl, trifluoromethyl and C1-6-alkoxy;

 $R^2$  is selected from hydrogen, halogen, optionally substituted aryl, optionally substituted aryloxy, and optionally substituted heteroaryl;

 $R^3$  is selected from hydrogen, optionally substituted  $C_{1.6}$ -alkoxy, halogen, cyano, and optionally substituted aryl, optionally substituted aryloxy, optionally substituted heteroaryl, amino,  $C_{1.6}$ -alkylcarbonylamino,  $C_{1.6}$ -alkylcarbonylamino, and mono- and di( $C_{1.6}$ -alkyl)aminosulfonyl; and

# with the proviso that R1, R2 and R3 are not all hydrogen;

X¹ and X² are independently selected from <u>hydroxy</u> (-OH) and acetoxy (-OAc) halogen, OR<sup>6</sup>;
OCOR<sup>5</sup>, N(R<sup>6</sup>)<sub>2</sub>, NHCOR<sup>5</sup>, NHSO<sub>2</sub>R<sup>5</sup>, and NHCON(R<sup>6</sup>)<sub>2</sub>, wherein R<sup>5</sup> is selected from C<sub>1-6</sub>-alkyl, optionally substituted aryl and optionally substituted heteroaryl, and each R<sup>6</sup> independently is selected from hydrogen, C<sub>1-6</sub>-alkyl, optionally substituted aryl and optionally substituted heteroaryl; and

pharmaceutically acceptable salts and prodrugs thereof.

Claim 25 (Currently amended): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IIb)

wherein

 $R^1$ ,  $R^2$ , and  $R^3$  independently are selected from hydrogen, optionally substituted  $C_{1:6}$ -alkyl, optionally substituted  $C_{2:6}$ -alkenyl, hydroxy, optionally substituted  $C_{1:6}$ -alkoxy, optionally substituted  $C_{1:6}$ -alkoxycarbonyl, optionally substituted  $C_{1:6}$ -alkoxycarbonyl, optionally substituted  $C_{1:6}$ -alkylcarbonyloxy, formyl, amino, monoand  $di(C_{1:6}$ -alkyl)amino, carbamoyl, mono- and  $di(C_{1:6}$ -alkyl)amino, carbamoyl, mono- and  $di(C_{1:6}$ -alkylsulphonylamino, cyano, carbamido, mono- and  $di(C_{1:6}$ -alkyl)-aminocarbonylamino,  $C_{1:6}$ -alkylsulphonylamino, cyano, carbamido, mono- and  $di(C_{1:6}$ -alkyl)-aminocarbonylamino,  $C_{1:6}$ -alkanoyloxy,  $C_{1:6}$ -alkylsulphonyl,  $C_{1:6}$ -alkylsulphinyl, aminosulfonyl, mono- and  $di(C_{1:6}$ -alkyl)-aminosulfonyl, nitro, optionally substituted  $C_{1:6}$ -alkylthio, and halogen, where any  $C_{1:6}$ -alkyl as an amino substituent is optionally substituted with hydroxy,  $C_{1:6}$ -alkoxy,

Application No. 10/599,121 Amendment dated December 29, 2009

Reply to Office Action of July 10, 2009

amino, mono- and di(C<sub>1-6</sub>-alkyl)amino, carboxy, C<sub>1-6</sub>-alkylcarbonylamino, C<sub>1-6</sub>-alkylaminocarbonyl, or halogen(s); and

or wherein  $R^1$  and  $R^2$  together with the carbon atoms to which they are attached form a heterocyclic ring, a heteroaromatic ring, an aromatic ring or a carbocyclic ring; and

# with the proviso that R1, R2 and R3 are not all hydrogen;

 $X^1$  and  $X^2$  are independently selected from  $\underline{hydroxy(-OH)}$  and  $\underline{acetoxy(-OAC)}$   $\underline{halogen, OR^6}$ ,  $OCOR^5$ ,  $N(R^6)_{a}$ ,  $NHCOR^5$ ,  $NHSO_2R^5$ ,  $\underline{and.NHCON(R^6)_{a}}$ ,  $\underline{wherein.R^6}$  is selected from  $C_{1-6}$ -alkyl, optionally substituted  $\underline{aryl.and.optionally.substituted.eryl.and.optionally.eryl.a$ 

pharmaceutically acceptable salts and prodrugs thereof.

Claim 26 (Withdrawn): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IIc)

wherein

 $R^{\perp}$  is selected from hydrogen, halogen,  $C_{1\text{-}6}$ -alkyl, trifluoromethyl and  $C_{1\text{-}6}$ -alkoxy;

 $R^2$  is selected from hydrogen, halogen, optionally substituted aryl, optionally substituted aryloxy, and optionally substituted heteroaryl;

Docket No.: 20517/0205421-US0

R<sup>3</sup> is selected from hydrogen, optionally substituted C<sub>1-6</sub>-alkoxy, halogen, cyano, and optionally substituted aryl, optionally substituted aryloxy, optionally substituted heteroaryl, amino, C<sub>1.6</sub>-alkylcarbonylamino, C1-6-alkylsulphonylamino, and mono- and di(C1-6-alkyl)aminosulfonyl; and

one of X1 and X2 is selected from halogen, OR6, OCOR5, N(R6), NHCOR5, NHSO<sub>2</sub>R5, and NHCON(R<sup>6</sup>)<sub>2</sub>, wherein R<sup>5</sup> is selected from C<sub>1-6</sub>-alkyl, optionally substituted aryl and optionally substituted heteroaryl, and each R<sup>6</sup> independently is selected from hydrogen, C<sub>1-6</sub>-alkyl, optionally substituted aryl and optionally substituted heteroaryl; and the other of X1 and X2 is selected from optionally substituted C1-6-alkyl, optionally substituted C2-6-alkenyl, carboxy, optionally substituted C1-6-alkoxycarbonyl, optionally substituted C1-6-alkylcarbonyl, formyl, carbamoyl, mono- and di(C1.6-alkyl)aminocarbonyl, cvano, aryl, arylcarbonyl, heterocyclyl, heterocyclylcarbonyl, heteroaryl, heteroarylcarbonyl, where any C1.6-alkyl as an amino substituent is optionally substituted with hydroxy, C1.6-alkoxy, amino, mono- and di(C1-6-alkyl)amino, carboxy, C1-6-alkylcarbonylamino, C1-6-alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted; and

pharmaceutically acceptable salts and prodrugs thereof.

Claim 27 (Withdrawn): A method of treating a mammal suffering from or being susceptible to cancer, the method comprising administering to the mammal a therapeutically effective amount of a 3,3-diphenyl-1,3-dihydro-indol-2-one type compound of the formula (IId)

wherein

R1. R2, and R3 independently are selected from hydrogen, optionally substituted C1.6-alkyl, optionally substituted C2-6-alkenyl, hydroxy, optionally substituted C1-6-alkoxy, optionally

substituted  $C_{2\cdot6}$ -alkenyloxy, carboxy, optionally substituted  $C_{1\cdot6}$ -alkoxycarbonyl, optionally substituted  $C_{1\cdot6}$ -alkylcarbonyloxy, formyl, amino, monoand di( $C_{1\cdot6}$ -alkyl)amino, carbamoyl, mono- and di( $C_{1\cdot6}$ -alkyl)aminocarbonyl,  $C_{1\cdot6}$ -alkylcarbonylamino,  $C_{1\cdot6}$ -alkylsulphonylamino, cyano, carbamido, mono- and di( $C_{1\cdot6}$ -alkyl)aminocarbonylamino,  $C_{1\cdot6}$ -alkylsulphonylamino, cyano, carbamido, mono- and di( $C_{1\cdot6}$ -alkyl)aminocarbonylamino,  $C_{1\cdot6}$ -alkanoyloxy,  $C_{1\cdot6}$ -alkylsulphonyl,  $C_{1\cdot6}$ -alkylsulphinyl, aminosulfonyl, mono- and di( $C_{1\cdot6}$ -alkyl)aminosulfonyl, nitro, optionally substituted  $C_{1\cdot6}$ -alkylthio, and halogen, where any  $C_{1\cdot6}$ -alkyl as an amino substituent is optionally substituted with hydroxy,  $C_{1\cdot6}$ -alkoxy, amino, mono- and di( $C_{1\cdot6}$ -alkyl)amino, carboxy,  $C_{1\cdot6}$ -alkylcarbonylamino,  $C_{1\cdot6}$ -alkylaminocarbonyl, or halogen(s); and

or wherein  $R^1$  and  $R^2$  together with the carbon atoms to which they are attached form a heterocyclic ring, a heteroaromatic ring, an aromatic ring or a carbocyclic ring; and

one of  $X^1$  and  $X^2$  is selected from halogen,  $OR^6$ ,  $OCOR^5$ ,  $N(R^6)_2$ ,  $NHCOR^5$ ,  $NHSO_2R^5$ , and  $NHCON(R^6)_2$ , wherein  $R^5$  is selected from  $C_{1\cdot6}$ -alkyl, optionally substituted aryl and optionally substituted heteroaryl, and each  $R^6$  independently is selected from hydrogen,  $C_{1\cdot6}$ -alkyl, optionally substituted aryl and optionally substituted heteroaryl; and the other of  $X^1$  and  $X^2$  is selected from optionally substituted  $C_{1\cdot6}$ -alkyl, optionally substituted  $C_{2\cdot6}$ -alkenyl, carboxy, optionally substituted  $C_{1\cdot6}$ -alkoxycarbonyl, optionally substituted  $C_{1\cdot6}$ -alkylcarbonyl, formyl, carboxyl, mono- and di( $C_{1\cdot6}$ -alkyl)aminocarbonyl, cyano, aryl, arylcarbonyl, heterocyclyl, heterocyclylcarbonyl, heteroarylcarbonyl, where any  $C_{1\cdot6}$ -alkyl as an amino substitutent is optionally substituted with hydroxy,  $C_{1\cdot6}$ -alkoxy, amino, mono- and di( $C_{1\cdot6}$ -alkyl)amino, carboxy,  $C_{1\cdot6}$ -alkyl-carbonylamino,  $C_{1\cdot6}$ -alkylaminocarbonyl, or halogen(s), and wherein any aryl, heterocyclyl and heteroaryl may be optionally substituted; and

pharmaceutically acceptable salts and prodrugs thereof.

Claim 28 (Currently amended): The method according to claim 1, wherein the compound is selected from Items 1 to 225 listed below:

- 5-Amino-6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 2 5-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

- 5-Fluoro-3.3-bis-(4-hydroxy-phenyl)-1.3-dihydro-indol-2-one:
- 4 3,3-Bis-(4-hydroxy-phenyl)-5-nitro-1,3-dihydro-indol-2-one;
- 5 6-Bromo-3.3-bis-(4-hydroxy-phenyl)-5.7-dimethyl-1.3-dihydro-indol-2-one:
- 6 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile;
  - 7 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-7-methyl-1,3-dihydro-indol-2-one;
  - 8 6-Bromo-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one:
    - 9 6-Bromo-5-ethyl-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-onc;
    - 10 6-Bromo-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-

# carbonitrile;

- 11 6-Bromo-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;
- 12 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;
- 13 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile:
  - 14 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-7-methyl-1,3-dihydro-indol-2-one;
  - 15 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;
  - 6-Chloro-5-ethyl-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 17 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile:
  - 18 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;
  - 19 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5-methyl-7-methoxy-1,3-dihydro-indol-2-one:
- 20 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-2-oxo-2,3-dihydro-1H-indole-5-carbonitrile:
  - 21 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-5-methyl-1,3-dihydro-indol-2-one;
  - 22 6-Chloro-5-ethyl-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-indol-2-one;
  - 23 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-5,7-dimethoxy-1,3-dihydro-indol-2-one;
  - 24 \_\_\_\_3,3-Bis-(4-hydroxy-phenyl)-1,3-dihydro-benzo[g]indol-2-one;
- 25 Acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-benzo[g|indol-3-yl]-phenyl ester;
  - 26 1-Amino-6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

- 27 Acetic acid 4-[3-(4-acetoxy-phenyl)-1-amino-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-vl]-phenyl ester:
- 28 Acetic acid 4-[3-(4-acetoxy-phenyl)-1-acetylamino-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-vll-phenyl ester:
  - 29 6-Chloro-7-cyclopropyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
  - 30 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-trifluoromethyl-1,3-dihydro-indol-2-one;
  - 31 6-Chloro-7-cyclopropoxy-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 32 6-(4-Fluoro-phenoxy)-3,3-bis-(4-hydroxy-phenyl)-7-trifluoromethyl-1,3-dihydroindol-2-one;
- 33 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-cyclopropyl-2-oxo-2,3-dihydro-1H-indol-3-yll-phenyl ester;
- 34 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-2-oxo-7-trifluoromethyl-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
- 35 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-cyclopropoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
- 36 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-(4-fluoro-phenoxy)-2-oxo-7-trifluoromethyl-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
  - 37 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-trifluoromethoxy-1,3-dihydro-indol-2-one;
- 38 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-2-oxo-7-trifluoromethoxy-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
  - 39 6-Chloro-4-fluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 40 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-4-fluoro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
- 41 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-4,7-dimethyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
- 42 6-Chloro-4,5-difluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 43 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-4,5-difluoro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
  - 44 3,3-Bis-(4-hydroxy-phenyl)-1,3,6,7,8,9-hexahydro-benzo[g]indol-2-one;

- 45 3.3-Bis-(4-hvdroxy-phenyl)-7-trifluoromethyl-1.3-dihydro-indol-2-one:
- 46 7-Chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 47 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-7-carbonitrile;
- 48 7-Ethyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 49 3,3-Bis-(4-hydroxy-phenyl)-7-morpholin-4-yl-1,3-dihydro-indol-2-one;
  - 50 3,3-Bis-(4-hydroxy-phenyl)-7-isopropyl-1,3-dihydro-indol-2-one;
- 51 7-tert-Butyl-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 52 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-7-carboxylic acid dimethylamide:
  - 53 3,3-Bis-(4-hydroxy-phenyl)-7-(4-methyl-piperazine-1-carbonyl)-1,3-dihydro-indol-

### 2-one;

- 54 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-carboxylic acid;
- 55 3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-1H-indole-5-carboxylic acid dimethylamide:
  - 56 3,3-Bis-(4-hydroxy-phenyl)-5-(morpholine-4-carbonyl)-1,3-dihydro-indol-2-one;
  - 57 3,3-Bis-(4-hydroxy-phenyl)-4-methoxy-1,3-dihydro-indol-2-one;
  - 58 3,3-Bis-(4-hydroxy-phenyl)-6-methoxy-1,3-dihydro-indol-2-one;
  - 59 3,3-Bis-(4-hydroxy-phenyl)-5-(4-methyl-piperazine-1-carbonyl)-1,3-dihydro-indol-

### 2-one;

- 60 3,3-Bis-(4-hydroxy-phenyl)-7-pyridin-3-yl-1,3-dihydro-indol-2-one;
- 61 7-Bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 62 7-Ethyl-5-fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 63 3,3-Bis-(4-hydroxy-phenyl)-5-iodo-1,3-dihydro-indol-2-one;
- 64 5-Amino-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 65 \_\_\_\_5-Amino-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 66 6-Bromo-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 67 7-Fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 68 3,3-Bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-indol-2-one;
- 69 4,7-Dichloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 70 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-1,7-dimethyl-1,3-dihydro-indol-2-one;

- 6-Chloro-3,3-bis-(4-fluoro-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 2 3,3-Bis-(4-hydroxy-phenyl)-7-(morpholine-4-carbonyl)-1,3-dihydro-indol-2-one;
- 73 3,3-Bis-(4-hydroxy-phenyl)-4,7-dimethyl-1,3-dihydro-indol-2-one;
- 74 3,3-Bis-(4-hydroxy-phenyl)-7-iodo-1,3-dihydro-indol-2-one;
- 75 3,3-Bis-(4-hydroxy-phenyl)-7-pyridin-4-yl-1,3-dihydro-indol-2-one;
- 76 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester;
  - 77 3,3-Bis-(4-hydroxy-phenyl)-5-phenyl-1,3-dihydro-indol-2-one;
  - 78 3,3-Bis-(4-hydroxy-phenyl)-7-thiophen-2-yl-1,3-dihydro-indol-2-one;
  - 79 3.3-Bis-(4-hydroxy-phenyl)-5-pyridin-4-yl-1.3-dihydro-indol-2-one:
  - 80 3.3-Bis-(4-hydroxy-phenyl)-5-thiophen-2-yl-1,3-dihydro-indol-2-one;
  - 81 5,7-Difluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
  - 82 6-Fluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
  - 83 3,3-Bis-(4-hydroxy-phenyl)-6-methoxy-7-methyl-1,3-dihydro-indol-2-one;
  - 84 6,7-Difluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
  - 85 6-Chloro-7-fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
  - 86 5-Fluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
  - 87 3,3-Bis-(4-hydroxy-phenyl)-5-methoxy-7-methyl-1,3-dihydro-indol-2-one;
  - 88 7-Chloro-3,3-bis-(4-hydroxy-phenyl)-4-methoxy-1,3-dihydro-indol-2-one;
  - 89 6-Fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
  - 90 N-[3,3-Bis-(4-hydroxy-phenyl)-2-oxo-2,3-dihydro-indol-1-yl]-acetamide;
- 91 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-6-yloxy]pentanoic acid methyl ester;
- 92 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-6-yloxy]-pentanoic acid;
- 93 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-5-yloxy]pentanoic acid methyl ester;
- 94 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-1H-indol-5-yloxy]pentanoic acid; and
  - 95 7-Chloro-6-fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one.

```
5 Amino 6 chloro 3,3 bis (4 hydroxy-phenyl) 7-methyl 1,3 dihydro indol-2-one
      5 Chloro 3,3 bis (4 hydroxy-phenyl) 7 methyl 1,3 dihydro indol-2-one
      5 Fluoro 3,3 bis (4 hydroxy-phenyl)-1,3-dihydro-indol-2-one
      4 3,3-Bis-(4-hydroxy-phenyl)-5-nitro-1,3-dihydro-indol-2-one
      5 3.3 Bis (4 hydroxy phenyl) 7 methyl 1.3 dihydro pyrrolo[3,2-clpyridin-2-one
      6 Bromo 3.3-bis (4-hydroxy-phenyl)-1.3-dihydro-pyrrolo[3,2-c]pyridin-2-one
      7 6 Bromo 3,3 bis (4 hydroxy phenyl)-7-methyl-1,3-dihydro-pyrrolo[3,2-e]pyridin-2-
           6 Bromo 3.3 bis (4-hydroxy-phenyl) 5.7-dimethyl-1.3-dihydro-indol-2-one
      9 6-Bromo-3,3-bis-(4-hydroxy-phenyl) 7-methyl-2-oxo-2,3-dihydro-1H-indole-5-
carbonitrile
             6 Bromo-3,3-bis-(4-hydroxy-phenyl) 5 methoxy 7 methyl-1,3-dihydro-indol-2-one
            6-Bromo-3,3-bis-(4-hydroxy-phenyl)-7-methoxy-1,3-dihydro-pyrrolo[3,2-e]pyridin-
      12 6 Bromo 7 ethyl 3.3 bis (4 hydroxy-phenyl) 1.3-dihydro-pyrrolo[3.2 elpyridin-2-
            6 Bromo 7 ethyl 3.3 bis (4 hydroxy-phenyl)-5-methyl 1.3 dihydro indol-2-one
      14 6 Bromo-5-ethyl-3.3-bis-(4-hydroxy-phenyl) 7-methyl-1.3-dihydro-indol 2-one
      15 6-Bromo 7-ethyl-3.3-bis (4-hydroxy-phenyl)-2-oxo-2.3-dihydro-1H-indole-5-
```

# carbonitrile

- 16 --- 6 Bromo 7 ethyl-3,3 bis (4 hydroxy phenyl) 5 methoxy-1,3 dihydro indol 2 one
- 17 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-e]pyridin-2-one
- 18 6-Chloro 3.3 bis (4-hydroxy-phenyl) 7 methyl 1.3-dihydro-pyrrolof 3.2-elpyridin 2-

### one

one

2-one:

one

- 19 6-Chloro-3,3-bis-(4-hydroxy-phenyl) 5,7 dimethyl-1,3-dihydro-indol-2-one
- 20 6 Chloro 3,3 bis (4 hydroxy phenyl) 7-methyl-2-oxo-2,3-dihydro-1H-indole-5carbonitrile
  - 21 6 Chloro 3.3 bis (4 hydroxy phenyl) 5-methoxy 7-methyl 1.3 dihydro-indol-2-one

- 22 6 Chloro 3,3-bis-(4-hydroxy-phenyl) 7 methoxy-1,3-dihydro-pyrrolo[3,2-e]pyridin-2-one
- 23 6-Chloro-7-ethyl-3,3-bis-(4-hydroxy-phenyl) 1,3-dihydro-pyrrolo[3,2-e]pyridin-2-one
  - 24 6 Chloro 7 ethyl-3,3 bis (4 hydroxy phenyl) 5 methyl-1,3 dihydro-indol 2 one
  - 25 6 Chloro-5-ethyl-3,3-bis (4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one
- 26 -- 6 Chloro 7 ethyl 3,3-bis (4-hydroxy-phenyl) 2 oxo 2,3 dihydro 1H-indole-5-

### carbonitrile

- 27 6 Chloro 7-ethyl 3,3-bis (4-hydroxy-phenyl)-5-methoxy 1,3-dihydro-indol-2-one
- 28 6-Chloro-3,3-bis (4-hydroxy-phenyl)-5-methyl-7-methoxy-1,3-dihydro-indol-2-one;
- 29 6 Chlore 3,3 bis (4 hydroxy phenyl) 7-methoxy 2-oxo 2,3-dihydro 1H indole 5-

## carbonitrile;

- 30 6-Chloro-3,3-bis (4-hydroxy-phenyl) 7-methoxy-1,3-dihydro-pyrrolo[3,2-o]pyridin-2-one:
  - 31 6-Chloro 3,3-bis (4-hydroxy-phenyl)-7-methoxy-5-methyl-1,3-dihydro-indol-2-one;
  - 32 6-Chloro-5-ethyl-3,3-bis (4-hydroxy-phenyl) 7-methoxy-1,3-dihydro-indol-2-one;
  - 33 6 Chloro 3.3 bis (4 hydroxy phenyl)-5.7-dimethoxy-1.3 dihydro indol 2-one:
- 34 N [4 [3 (4 Acetylamino-phenyl) 5-chloro 7-methyl 2 oxo 2,3-dihydro-1H-indol-3-yll-phenyl}-acetamide:
- 35 N [4-[5-Chloro-3 (4-methanesulfonylamino-phenyl) 7-methyl-2-oxo-2,3-dihydro-HH-indol-3-yl]-phenyl]-methanesulfonamide
- 36 N-{4-{3-(4-Acetylamino-phenyl)-6-chloro-7-methyl-2-oxo-2;3-dihydro-1H-indol-3yll-phenyll-acetamide:
- 37 N [4-[6-Chloro-3-(4-methanesulfonylamino-phenyl) 7-methyl-2-oxo-2,3-dihydro-HH-indol-3-yl|-phenyl}-methanesulfonamide;
- 38 N [4-[3 (4-Acetylamino-phenyl) 5-chloro-7-methoxy-2-oxo-2,3-dihydro-1H-indol-3-yll-phenyl)-acetamide:
- 39 N-{4-[5-Chloro-3-(4 methanesulfonylamino-phenyl)-7-methoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-methanesulfonamide;

- 40 N [4 [3 (4-Acetylamino-phenyl) 6-chloro-7-methoxy-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide; and
- 41 N. (4-[6-Chloro-3 (4 methanesulfonylamino-phenyl)-7-methoxy-2-oxo-2,3-dihydro-1H indol-3-yl]-phenyl]-methanesulfonamide
- 42——2-Chloro-6,6-bis-(4-hydroxy-phenyl) 3-methyl-4,6-dihydro-3H-pyrrolo[2,3-dlimidazol-5-one
- 43 Acetic acid 4 [6 (4 acetoxy-phenyl) 2 chloro-3-methyl-5-oxo-3,4,5,6-tetrahydropyrrolo[2,3-d]imidazol-6-yl]-phenyl-ester
- 44 6,6-Bis (4-amino-phenyl) 2-chloro-3-methyl-4,6-dihydro-3H-pyrrolo[2,3-dlimidazol-5-one
- 45 2-Chloro-6,6-bis-(4-dimethylamino-phenyl)-3-methyl-4,6-dihydro-3H-pyrrolo[2,3-dlimidazol-5-one
- 46 N-{4-[6-(4-Acetylamino-phenyl)-2-chloro-3-methyl-5-oxo-3,4,5,6-tetrahydro-pyrrolo[2,3-d]imidazol-6-yl]-phenyl}-acetamide
- 47 N [4 [2-Chloro-6 (4-methanesulfonylamino-phenyl)-3-methyl-5-oxo-3,4,5,6-tetrahydro-pyrrolo[2,3-d]imidazol-6-yl]-phenyl]-methanesulfonamide
  - 48 4,4 Bis (4 hydroxy-phenyl)-1-methyl 4,6-dihydro-1H-pyrrolo[2,3-e]pyrazol-5-one
- 49 Acetic acid 4-[4 (4-acetoxy-phenyl) 1-methyl-5-oxo-1,4,5,6-tetrahydro-pyrrolo[2,3-elovrazol-4-vll-phenyl-ester
  - 50 4,4 Bis (4-amino-phenyl) 1-methyl-4,6-dihydro-HI-pyrrolo[2,3-c]pyrazol-5-one
- 51 N-{4-[4-(4 Methanesulfonylamino-phenyl) 1 methyl-5-oxo-1,4,5,6-tetrahydro-pyrrolo[2,3-clpyrazol 4 yll-phenyl) methanesulfonamide
- 52 - 4,4 Bis (4 dimethylamino-phenyl) 1 methyl 4,6 dihydro-1H-pyrrolo[2,3-e]pyrazol-5-one
- 53 N [4 [4 (4 Acetylamino-phenyl) 1-methyl-5-oxo-1,4,5,6-tetrahydro-pyrrolo[2;3-e]pyrazol-4-yl] phenyl] acetamide
  - 54 4,4-Bis-(4-hydroxy-phenyl) 2-methyl-2,6-dihydro-4H-pyrrolo[2,3-c]pyrazol-5-one
- 55 Acetic-acid 4-[4 (4-acetoxy-phenyl) 2-methyl-5-oxo 2,4,5,6-tetrahydro-pyrrolo[2,3-e]pyrazol-4-yl]-phenyl-ester
  - 56 4,4 Bis (4 amino-phenyl) 2-methyl 2,6-dihydro-4H-pyrrolo[2,3-c]pyrazol-5-one

- Application No. 10/599,121 Docket No.: 20517/0205421-US0
- 57 4,4-Bis (4-dimethylamino-phenyl)-2-methyl-2,6-dihydro-4H-pyrrolo[2,3-o]pyrazol-5-one
- 58 N-[4-[4-(4-Acetylamino-phenyl)-2-methyl-5-oxo-2,4,5,6-tetrahydro-pyrrolo[2,3-elpyrazol-4-yl]-phenyl]-acetamide
- 59 N {4 [4 (4 Methanesulfonylamino-phenyl) 2-methyl-5-oxo-2,4,5,6-tetrahydro-pyrrolo[2,3-e]pyrazol-4-yl]-phenyl]-methanesulfonamide
  - 60 4,4 Bis (4-hydroxy-phenyl)-4,6-dihydro-thieno[2,3-b]pyrrol-5-one
- 61—Acetic acid 4 [4 (4-acetoxy-phenyl)-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yll-phenyl-ester
  - 62 4,4-Bis-(4-amino-phenyl)-4,6-dihydro-thieno[2,3-b]pyrrol-5-one
  - 63 4,4-Bis-(4-dimethylamino-phenyl) 4,6-dihydro-thieno[2,3-b]pyrrol-5-one
- 64 N-{4-[4-(4-Acetylamino-phenyl)-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]phenyl)-acetamide
- 65 N {4 [4-(4-Methanesulfonylamino-phenyl)-5-exo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl}-methanesulfonamide
  - 66 2 Chloro 4.4 bis (4-hydroxy-phenyl) 4.6-dihydro-thieno[2.3-b]pyrrol-5-one
- 67 Acetic acid 4-[4-(4-acetoxy-phenyl) 2-chloro-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl ester
  - 68 4,4-Bis-(4-amino-phenyl)-2-chloro-4,6-dihydro-thieno[2,3-b]pyrrol-5-one
  - $\textcolor{red}{69} \textcolor{red}{2-Chloro-4,4-bis-(4-dimethylamino-phenyl)-4,6-dihydro-thiono} \textcolor{blue}{[2,3-b]pyrrol-5-one}$
- 70 N-{4-{4-(4-Acetylamino-phenyl) 2-chloro-5-oxo-5,6-dihydro-4H-thieno[2,3-b|pyrrol-4-yll-phenyl}-acetamide
- $71 N (4-[2-Chloro-4-(4-methanesulfonylamino-phenyl)-5-oxo-5,6-dihydro-4H-thieno[2,3-b]pyrrol-4-yl]-phenyl}-methanesulfonamide$ 
  - 72 4,4 Bis (4-hydroxy-phenyl) 4,6 dihydro-furo[2,3-b]pyrrol-5-one
- 73 Acetic acid 4-[4-(4-acetoxy-phenyl)-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yl]phenyl-ester
  - 74 4,4 Bis (4 amino-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one
  - 75 4,4-Bis (4 dimethylamino-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one

phenyl)-acetamide

- 76 N. (4-[4-(4-Acetylamino-phenyl)-5-oxo-5.6-dihydro-4H-furo[2,3-bloyrrol-4-yl]-
- 77 N (4-[4-(4-Methanesulfonylamino-phenyl)-5-oxo-5,6-dihydro-4H furo[2,3-b]pyrrol-4-yl]-phenyl}-methanesulfonamide
  - 78 2 Chloro-4,4-bis-(4-hydroxy-phenyl)-4,6-dihydro-furo[2,3-b]pyrrol-5-one
- 79 Acetic acid 4-[4-(4-acetoxy-phenyl)-2-chloro-5-oxo-5,6-dihydro-4H-furo[2,3-blovrol-4-vl]-phenyl-ester
  - 80 4.4-Bis-(4-amino-phenyl) 2-chloro-4.6-dihydro-furo[2,3-blpyrrol-5-one
  - 81 2 Chloro 4.4 bis (4-dimethylamino-phenyl) 4,6-dihydro-furo[2,3-b]pyrrol-5-one
- 82 N (4 [4 (4 Acetylamino-phenyl) 2-chloro-5-oxo-5,6-dihydro-4H-furo[2,3-b]pyrrol-4-yll-phenyl)-acetamide
- 83 N. [4 [2 Chlore 4 (4-methanesulfonylamino-phenyl) 5-oxe-5,6-dihydre-4Hfuro[2,3-blpyrrol-4-yl]-phenyl)-methanesulfonamide
  - 84 3,3 Bis (4 hydroxy phenyl) 6-methyl 3,8 dihydro 1H 1,8 diaza-as-indacen 2-one
- 85 Acetic acid 4-[3-(4-acetoxy-phenyl) 6 methyl 2-exo-1,2,3,8 tetrahydro-1,8-diaza-asindacen-3-vll-phenyl ester
  - 86 3,3-Bis (4-amino-phenyl) 6-methyl-3,8-dihydro-1H-1,8-diaza-as-indacen-2-one
- 87 3,3-Bis-(4-dimethylamino-phenyl) 6-methyl-3,8-dihydro-1H-1,8-diaza-as-indacen-2-one
- 88 N- [4-[3-(4-Acetylamino-phenyl)-6-methyl-2-oxo-1,2,3,8-tetrahydro-1,8-diaza-as-indacen-3-yl]-phenyl}-acetamide
- $89 N \{4 \ [3 \ (4 \ Methanesulfonylamino-phenyl) \ 6 \ methyl \ 2 \ oxo \ 1,2,3,8 \ tetrahydro-1,8-diaza-as-indacen \ 3-yl]-phenyl} methanesulfonamide$ 
  - 90 --- 3,3-Bis-(4-hydroxy-phenyl) 1,3-dihydro-benzo[g]indol-2-one
- $91 Acetie acid 4 \left[3 (4 acetoxy-phenyl) 2 oxo 2, 3 dihydro 1H benzo [g]indol 3 yl] phenyl ester$ 
  - 92 3,3-Bis-(4-amino-phenyl)-1,3-dihydro-benzo[g]indol-2-one
  - 93 3,3-Bis-(4-dimethylamino-phenyl) 1,3-dihydro-benzo[g]indol-2-one
- 94 N-{4-{3-(4-Acetylamino-phenyl})-2-oxo-2,3-dihydro-1H-benzo{g}indol-3-yl}-phenyl}-acetamide

- 95 N. [4 [3 (4-Methanesulfonylamino-phenyl)-2-oxo-2,3-dihydro-1H-benzo[g]indol-3-yl]-phenyl}-methanesulfonamide
  - 96 1-Amino-6-chloro 3.3-bis (4 hydroxy-phenyl) 7-methyl-1,3-dihydro-indol-2-one
- 97 Acetic acid 4-[3-(4-acetoxy phenyl) 1 amino 6-chloro 7 methyl 2 oxo 2,3-dihydro-HH indol-3 vll-phenyl ester
- 98 N (4 [3 (4-Acetylamino-phonyl) 1 amino 6 chloro-7-methyl 2 oxo 2,3 dihydro-1H-indol-3-yll-phonyl) acetamide
- 99 N-{4-[1 Amino 6 chloro 3 (4 methanesulfonylamino phenyl) 7-methyl 2-oxo 2,3-dihydro-1H indol 3-yl}-phenyl}-methanesulfonamide
- 100 Acetic acid 4 [3 (4 acetoxy-phenyl)-1-acetylamine-6-chlore-7-methyl-2-oxo-2,3-dihydro-1H indol-3-yll-phenyl-ester
- 101 N [3,3-Bis (4-amino-phenyl) 6-ehloro-7-methyl-2-oxo-2,3-dihydro-indol-1-yl]-acetamide
- 102 N [6 Chloro 3,3-bis (4 dimethylamino-phenyl) 7-methyl 2-oxo 2,3 dihydro-indol 1-yll-acetamide
- 103 N [3,3-Bis (4-acetylamino-phenyl)-6-chloro 7 methyl 2-oxo-2,3-dihydro-indol-1-yll-acetamide
- 104 N [6 Chlore-3,3-bis (4 methanesulfonylamino-phenyl) 7 methyl 2 exe 2,3 dihydro-indol 1 yll-acetamide
  - 105 6-Chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indole-2-thione
- 106 Acetic acid 4 [3 (4-acetoxy-phenyl) 6-chloro-7-methyl-2-thioxo-2,3-dihydro-1Hindol-3-yl]-phenyl-ester
  - 107 3,3 Bis (4-amino phenyl) 6 chloro 7 methyl 1,3-dihydro-indole 2-thione
  - $108 6 \cdot Chloro 3, 3 \cdot bis \cdot (4 \cdot dimethylamino \cdot phenyl) \cdot 7 \cdot methyl \cdot 1, 3 \cdot dihydro \cdot indole \cdot 2 \cdot thione$
- 109 N (4-[3 (4 Acetylamino phenyl) 6-chloro-7-methyl-2-thioxo-2,3-dihydro-1H-indol-3-yl]-phenyl}-acetamide
- 110 Methanesulfonic acid 4 [6-chloro-3 (4 methanesulfonyloxy-phenyl) 7 methyl-2-thicxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester
- 111 Acetic acid 4-[4-(4-acetoxy-phenyl) 2-chloro-5-thioxo-5,6-dihydro-4H-thieno[2,3-b|pyrrol-4-yl]-phenyl-ester

b|pyrrol-4-yl| phenyl ester

- 112 Acetic acid 4-[4-(4-acetoxy-phenyl) 2-chloro-5-thioxo-5,6-dihydro-4H-fure[2,3-
- 113 6,6-Bis (4-amino-phonyl) 2-chloro-3-methyl 4,6-dihydro-thieno[3,2-b]pyrrole-5-thiene
- 114 2 Chloro 6,6-bis-(4-dimethylamino-phenyl)-3-methyl-4,6-dihydro-3H-pyrrolo[2,3-dlimidazole-5-thione
- 115 N-{4-[6-(4-Acetylamino-phenyl)-3-chloro-5-thioxo-1,4,5,6-tetrahydro-pyrrolo[3,2-clpyrazol-6-yl]-phenyl}-acetamide
- 116 Methanesulfonie acid 4-[2-chlore 4-(4-methanesulfonyloxy-phenyl) 5-thioxe-5,6dihydro 4H-furo[2,3-b]pyrrol-4-yl]-phenyl-ester
  - 117 6-Chloro-7-cyclopropyl-3,3-bis (4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
- 118 6-Chloro-7-cyclopropyl-3,3-bis (4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-chvridin-2-one
  - 119 6-Chloro 3,3-bis (4-hydroxy-phenyl)-7-trifluoromethyl-1,3-dihydro-indol-2-one
- 120 6 Chloro-3,3-bis (4-hydroxy-phenyl) 7-trifluoromethyl-1,3-dihydro-pyrrolo[3,2-e|pyridin-2-one
  - 121 6-Chloro-7-eyelopropoxy-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
- 122 6-Chloro-7-eyelopropoxy-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-pyrrolo[3,2-elpyridin-2-one
- 123 6 (4 Fluoro-phenoxy) 3,3 bis (4 hydroxy-phenyl) 7-trifluoromethyl 1,3-dihydro-indol-2-one
- 124 Acetic acid 4 [3-(4-acetoxy-phenyl) 6-chloro 7-cyclopropyl-2-oxo 2,3-dihydro 1H-indol-3-yl] phenyl ester
- 125 Acetic acid 4 [3 (4 acetoxy-phenyl) 6 chloro 7-cyclopropyl 2-oxo-2,3-dihydro-1H-pyrrolo[3,2-e]pyridin 3-yl]-phenyl ester
- 126 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro 2-oxo 7-trifluoromethyl 2,3-dihydro-1H-indol-3-yl]-phenyl-ester
- 127 Acetic acid 4-[3-(4-acetoxy-phenyl)-6-chloro-2-oxo-7-trifluoromethyl-2,3-dihydro-1H-pyrrolo[3,2-e]pyridin-3-yl]-phenyl-ester

- 128 Acetic acid 4 [3-(4-acetoxy-phenyl)-6-chloro-7-cyclopropoxy-2-oxo-2,3-dihydro-1H-indol-3-vll-phenyl ester
- 129 Acetic acid 4 [3-(4-acetoxy-phenyl) 6-chloro-7-cyclopropoxy-2-oxo-2,3-dihydro-1H-pyrrolo[3,2-c]pyridin-3-yl]-phenyl-ester
- 130 Acetic acid 4 [3 (4 acetoxy-phenyl) 6 (4 fluoro-phenoxy)-2-oxo-7-trifluoromethyl-2-3-dihydro-1H-indol-3-vll-phenyl-ester
- 131 Dimethylamino-acetic acid 4 (6-chloro 7-cyclopropyl 3 [4 (2-dimethylamino-acetoxy) phenyl] 2-oxo-2,3-dihydro-1H-indol-3-yl} phenyl ester
- 132 Dimethylamino acetic acid 4 (6-chloro-7-cyclopropyl-3 [4 (2-dimethylamino-acetoxy)-phenyl]-2-oxo-2,3-dihydro-1H-pyrrolo[3,2-c]pyridin-3-yl}-phenyl-ester
- 133 Dimethylamino-acetic-acid 4 (6-chloro-3-[4 (2-dimethylamino-acetoxy) phenyl] 7-methyl-2-oxo-2,3-dihydro-HH-indol-3-yl} phenyl-ester
  - 134 6 Chloro 3,3 bis (4-hydroxy-phenyl)-7-trifluoromethoxy-1,3-dihydro-indol-2-one
- 135 Acetic acid 4 [3 (4 acetoxy-phenyl) 6 chloro-2-oxo 7-trifluoromethoxy 2,3-dihydro-1H-indol-3 yl] phenyl ester
- 136 Dimethylamino acetic acid 4 (6 chloro 3 [4 (2 dimethylamino acetoxy) phenyl] 2exe-7-triflueremethexy 2.3 dihydro 1H indol 3 vll-phenyl ester
  - 137 6-Chloro-4-fluoro 3,3 bis (4 hydroxy phenyl) 7-methyl-1,3 dihydro-indol 2-one
- 138 3 Chloro 7,7 bis (4 hydroxy-phenyl) 4-methyl-5,7-dihydro-pyrrolo[3,2-e]pyridazin-6-one
- $\frac{139 Acetic acid 4 \left[3 (4 acetoxy-phenyl) 6 chloro 4 fluoro 7 methyl 2 oxo 2,3 dihydro-1H indol 3 yl] phenyl ester}{}$
- 140 Acetic acid 4-[3-(4-acetoxy-phenyl) 6-chloro 4,7-dimethyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester
- $141 Acetic acid 4 \cite{T-(4-acetoxy-phenyl)-3-chloro-4-methyl-6-oxo-6,7-dihydro-5H-pyrrolo[3,2-e]pyridazin-7-yl]-phenyl-ester$ 
  - 142 6-Chloro-4,5-difluoro-3,3-bis (4-hydroxy phenyl)-7-methyl-1,3-dihydro-indol-2-one
- 143 Acetic acid 4-[3-(4 acetoxy-phenyl) 6-chloro 4,5-difluoro 7-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl-ester
  - 144 3,3-Bis-(4 hydroxy phenyl) 3,6,7,8 tetrahydro 1H 1 aza as indacen 2-one

- Docket No.: 20517/0205421-US0
- 145 3,3-Bis-(4-hydroxy-phenyl) 1,3,6,7,8,9-hexahydro-benzo[g]indol-2-one
- 146 3,3-Bis-(4-hydroxy phenyl)-7-trifluoromethyl-1,3-dihydro-indol-2-one
- 147 7-Chloro 3.3 bis (4-hydroxy-phenyl) 1.3 dihydro-indol 2 one
- 148 3.3-Bis-(4-hydroxy phenyl)-2-oxo-2,3-dihydro-1H-indole-7-carbonitrile
- 149 7-Ethyl-3,3 bis (4-hydroxy-phenyl)-1,3-dihydro-indol 2 one
- 150 3,3-Bis-(4-hydroxy-phenyl) 7-morpholin 4-yl-1,3-dihydro-indol-2-one
- 151 3,3-Bis (4-hydroxy phenyl)-7-isopropyl-1,3-dihydro-indol-2-one
- 152 7-tert-Butyl-3,3-bis (4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
- 153 3,3-Bis (4-hydroxy-phenyl) 2-oxo-2,3-dihydro-1H-indole-7-carboxylic acid dimethylamide
- 154 3,3-Bis (4-hydroxy-phenyl) 7 (4-methyl-piperazine 1-carbonyl) 1,3-dihydro-indol-2-one
  - 155 3.3-Bis (4 hydroxy-phenyl) 2-oxo-2.3-dihydro-1H-indole-5-carboxylic acid
- 156 3,3 Bis (4-hydroxy-phenyl) 2-oxo-2,3-dihydro 1H indole-5-earboxylio acid dimethylamide
  - 157 3.3 Bis-(4-hydroxy phenyl)-5-(morpholine-4-carbonyl)-1.3 dihydro-indol-2-one
  - 158 3.3 Bis (4-hydroxy-phenyl) 4-methoxy-1,3-dihydro-indol-2-one
  - 159 3,3-Bis-(4-hydroxy-phenyl)-6-methoxy-1,3-dihydro-indol-2-one
  - 160 3,3-Bis (4-hydroxy-phenyl) 5-(4-methyl-piperazine-1-carbonyl) 1,3-dihydro-indol-

#### 2-one

- 161 6-Chloro-3.3-bis-(4-mercapto-phenyl)-7-methyl-1.3-dihydro-indol-2-one
- 162 N {4 [3 (4 Acetylamino phenyl) 7 methyl 2-oxo 2,3 dihydro HH indol-3-yl]phenyl)-acetamide
  - 163 3.3-Bis-(4-hydroxy-phenyl) 7 (3-methoxy-prop-1-ynyl) 1.3-dihydro-indol-2-one
  - 164 3.3 Bis (4-hydroxy-phenyl)-7-pyridin-3-yl-1,3-dihydro-indol-2-one
  - 165 7-Bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
  - 166 6-Chloro-3,3-bis (4-methanesulfonyl-phenyl) 7-methyl-1,3-dihydro-indol-2-one
  - 167 6,6-Bis-(4-hydroxy phenyl)-4,6-dihydro-pyrrolo[3,2-d]thiazol-5-one
  - 168 6,6-Bis-(4-hydroxy-phenyl) 2-methyl-4,6-dihydro-pyrrolo[3,2-d]thiazol-5-one
  - 169 6,6 Bis-(4-hydroxy phenyl) 2-isopropyl-4,6-dihydro-pyrrolo[3,2-d]thiazol 5-one

170 2-Chloro 6.6 bis (4 hydroxy phenyl) 4.6 dihydro-pyrrolof 3.2 dlthiazol 5 one

Docket No : 20517/0205421-US0

- 171 4,4-Bis (4-hydroxy-phenyl) 4,6-dihydro-pyrrolo[3,2-d]isothiazol-5-one
- 172 3,3-Bis-(4-hydroxy-phenyl) 7-methyl-1,3-dihydro-pyrrolo[2,3-e]pyridin-2-one
- 173 3.3 Bis (4 hydroxy-phenyl) 7 methyl-1.3-dihydro-pyrrolo[3,2 b]pyridin-2-one
- 174 3.3-Bis (4 fluoro-phenyl) 7-methyl-1.3 dihydro-pyrrolo[3,2-b]pyridin-2-one
- 175 3.3-Bis (4-fluoro-phenyl) 7-methyl-1,3-dihydro-pyrrolo[3,2-c]pyridin 2-one
- 176 3,3 Bis (4-fluoro-phenyl) 7-isopropyl-1,3-dihydro-pyrrolo[3,2-c]pyridin-2-one
- 177 3,3 Bis (4 hydroxy-phenyl) 3,6,7,8 tetrahydro 1H-1,5 diaza as indacen 2 one
- 177 5,5-Bis (1-flydroxy-pherly)-5,0,7,5 tellarlydro 111 1,5 diaza as medden 2 one
- 178 3,3-Bis (4-hydroxy-phenyl)-3,6,7,8-tetrahydro-1H-1,4-diaza-as-indacen-2-one
- $179 3,3 Bis. \\ (4-hydroxy-phenyl) 1,3,6,7,8,9-hexahydro-pyrrolo[3,2-e] \\ quinolin-2-one$
- 180 3,3-Bis (4-hydroxy-phenyl)-1,3,6,7,8,9-hexahydro-pyrrolo[3,2-c]isoquinolin-2-one
- 181 5-Fluoro 3,3 bis (4 hydroxy-phenyl) 3,6,7,8-tetrahydro 1H-1 aza as indacen 2-one
- 182 7-Ethyl-5-fluoro-3.3 bis (4-hydroxy-phenyl)-1.3-dihydro-indol-2-one
- 183 3.3 Bis (4-hydroxy-phenyl)-1.3.6.8 tetrahydro-7-oxa 1-aza-as-indaeen-2-one
- 184 3.3-Bis (4-hydroxy-phenyl)-1.3.7.8-tetrahydro-6-oxa-1-aza-as-indacen-2-one
- 185 3,3 Bis (4-hydroxy-phenyl) 1,6,7,9 tetrahydro 3H 8 oxa-1-aza-

## eyelopenta[a]naphthalen-2-one

- 186 3.3 Bis (4-hydroxy-phenyl) 1.7.8.9 tetrahydro-3H pyrano[2,3-g]indol-2-one
- 187 3,3-Bis (4-hydroxy-phenyl) 7-methyl 3,6,7,8-tetrahydro-HI-1,7-diaza as indacen-2-one
- 188 3,3-Bis (4-hydroxy-phenyl) 7-methyl-1,3,7,8-tetrahydro 1,7-diaza-as-indacene 2,6-dione
- 189 3,3-Bis (4-hydroxy-phenyl) 7,8,8-trimethyl 1,3,7,8-tetrahydro 1,7-diaza-asindacene 2,6-dione
  - 190 3,3-Bis-(4-hydroxy-phenyl)-5-iodo-1,3-dihydro-indol-2-one
  - 191 5-Amino 3,3 bis (4 hydroxy-phenyl) 1,3 dihydro-indol 2-one
  - 192 5-Amino 3.3-bis (4-hydroxy phenyl)-7-methyl-1.3-dihydro indol-2-one
  - 193 6 Bromo 3,3 bis (4 hydroxy phenyl) 7 methyl-1,3 dihydro-indol-2 one
  - 194 7-Fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
  - 195 3,3 Bis (4 hydroxy-phenyl) 7-methoxy 1,3-dihydro-indol-2-one

- Docket No.: 20517/0205421-US0
  - 196 4.7 Dichloro 3.3 bis (4 hydroxy phenyl) 1.3-dihydro indol 2 one
  - 197 6 Chloro-3,3-bis-(4-hydroxy-phenyl) 1,7-dimethyl-1,3-dihydro-indol-2-one
  - 198 6 Chloro 3,3 bis (4 fluoro phenyl) 7 methyl-1,3-dihydro-indol-2 one
  - 199 3,3 Bis (4-hydroxy-phenyl)-7 (morpholine 4-carbonyl)-1,3-dihydro-indol-2-one
  - 200 3.3 Bis (4 hydroxy phenyl) 1.3 dihydro pyrrolo[2.3 dlpyridin 2 one
- 201 N (4 f6 Chloro 3 (4 methanesulfonylamino phenyl) 7 methyl 2 oxo 2,3 dihydro-1H-indol-3-vl]-phenvl)-methanesulfonamide
  - 202 3,3-Bis (4-hydroxy-phenyl) 4,7-dimethyl 1,3-dihydro-indol-2-one
  - 203 3,3-Bis-(4-hydroxy-phenyl)-7-iodo-1,3-dihydro-indol-2-one
  - 204 3.3-Bis-(4-hydroxy-phenyl)-7-pyridin-4-yl-1.3-dihydro-indol-2-one
- 205 Acetic acid 4 [3 (4 acetoxy-phenyl) 6-chloro 7-methyl 2-oxo 2,3 dihydro 1H-indol-3-vll-phenyl ester
  - 206 3.3 Bis (4 hydroxy phenyl) 5 phenyl-1.3 dihydro-indol-2 one
  - 207 3.3 Bis (4-hydroxy-phenyl) 7-thiophen 2-yl 1.3-dihydro-indol-2-one
  - 208 3.3-Bis-(4-hydroxy-phenyl)-5-pyridin-4-yl-1,3-dihydro-indol-2-one
  - 209 3,3 Bis-(4-hydroxy-phenyl) 5-thiophen 2-yl-1,3-dihydro-indol-2-one
  - 210 5,7-Difluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
  - 211 6-Fluoro-3.3-bis-(4-hydroxy-phenyl) 7-methyl 1.3-dihydro-indol-2-one
  - 212 3.3-Bis-(4-hydroxy-phenyl)-6-methoxy-7-methyl-1.3-dihydro-indol-2-one
  - 213 6.7-Difluoro-3.3-bis-(4-hydroxy-phenyl) 1.3-dihydro-indol-2-one
  - 214 6 Chloro 7 fluoro 3,3 bis (4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
  - 215 5-Fluoro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one
  - 216 3.3 Bis (4 hydroxy phenyl) 5 methoxy 7-methyl-1.3-dihydro-indol-2-one
  - 217 3.3 Bis (4 hydroxy phenyl) 1.3 dihydro pyrrolo[2,3 b]pyridin 2 one
  - 218 7-Chloro-3,3-bis-(4-hydroxy-phenyl) 4-methoxy-1,3-dihydro-indol-2-one
  - 219 6-Fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one
  - 220 N. [3,3-Bis-(4-hydroxy-phenyl) 2-oxo-2,3-dihydro-indol-1-yl]-acetamide
- 221 5-[3,3-Bis-(4-hydroxy-phenyl) 7-methyl 2-oxo 2,3-dihydro-1H-indol-6-yloxy]pentanoic acid methyl ester

23

Application No. 10/599,121 Docket No.: 20517/0205421-US0

222 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-HH-indol-6-yloxy]pentanoic acid

223 — 5-[3,3-Bis-(4-hydroxy-phenyl)-7-methyl-2-oxo-2,3-dihydro-HH-indol-5-yloxy]pentanoic-acid-methyl-ester

224 — 5 [3,3-Bis (4-hydroxy-phenyl) 7-methyl 2-oxo-2,3-dihydro-1H-indol-5-yloxy]pentanoic-acid

225 7-Chloro-6-fluoro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one.

Claim 29 (Currently amended): The method according to claim 1, wherein the <u>method</u> medicament further comprises <u>administering</u> one or more other chemotherapeutic agents.

Claim 30 (canceled).

Claim 31 (Withdrawn): A compound of the general formula (I)

(I)

as defined in claim 1, with the proviso that the compound is not one selected from

- 3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,
- 3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;
- 3.3-bis-(4-hvdroxy-phenyl)-4.5-dimethyl-1.3-dihydro-indol-2-one:
- 3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;
- 5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;
- 3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;
- 3.3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

Docket No.: 20517/0205421-US0

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one; acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 32 (Withdrawn): A 3.3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)

as defined in claim 24, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3.3-bis-(4-hydroxy-phenyl)-4.5-dimethyl-1.3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one; 3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 33 (Withdrawn): A pharmaceutical composition comprising a compound as defined in claim 1 and a pharmaceutically acceptable carrier.

Claim 34 (Withdrawn): A 3.3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)

Application No. 10/599,121 Docket No.: 20517/0205421-US0 Amendment dated December 29, 2009

Reply to Office Action of July 10, 2009

as defined in claim 25, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 35 (Withdrawn): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)

as defined in claim 26, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

Claim 36 (Withdrawn): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)

(II)

as defined in claim 27, with the proviso that the compound is not one selected from:

3, 3-b is-(4-hydroxy-phenyl)-1, 3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

acetic acid 4-[3-(4-acetoxy-phenyl)-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester; and acetic acid 4-[3-(4-acetoxy-phenyl)-5-methyl-2-oxo-2,3-dihydro-1H-indol-3-yl]-phenyl ester.

lication No. 10/599,121 Docket No.: 20517/0205421-US0

Claim 37 (Withdrawn): A 3,3-Diphenyl-1,3-dihydro-indol-2-one type compound of the formula (II)

as defined in claim 28, with the proviso that the compound is not one selected from:

3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one,

3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-4,5-dimethyl-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5,7-dimethyl-1,3-dihydro-indol-2-one;

5-bromo-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

5-chloro-3,3-bis-(4-hydroxy-phenyl)-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methoxy-1,3-dihydro-indol-2-one;

3,3-bis-(4-hydroxy-phenyl)-5-methyl-1,3-dihydro-indol-2-one;

6-chloro-3,3-bis-(4-hydroxy-phenyl)-7-methyl-1,3-dihydro-indol-2-one;

 $acetic\ acid\ 4\hbox{-}[3\hbox{-}(4\hbox{-}acetoxy\hbox{-}phenyl)\hbox{-}2\hbox{-}oxo\hbox{-}2,3\hbox{-}dihydro\hbox{-}1H\hbox{-}indol\hbox{-}3\hbox{-}yl]\hbox{-}phenyl\ ester; and$ 

 $acetic\ acid\ 4\hbox{-}[3\hbox{-}(4\hbox{-}acetoxy\hbox{-}phenyl)\hbox{-}5\hbox{-}methyl\hbox{-}2\hbox{-}oxo\hbox{-}2,3\hbox{-}dihydro\hbox{-}1H\hbox{-}indol\hbox{-}3\hbox{-}yl]\hbox{-}phenyl\ ester.$ 

Claim 38 (Currently amended): The method according to claim 1, wherein both of  $X^1$  and  $X^2$  are hydroxyl (-OH).

Claim 39 (New): The method according to claim 1, wherein R<sup>4</sup> is hydrogen.

Claim 40 (New): The method according to claim 39, wherein R<sup>3</sup> and R<sup>4</sup> are both hydrogen.